

Abhijit Mitra

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	On the Nature of Nucleobase Stacking in RNA: A Comprehensive Survey of Its Structural Variability and a Systematic Classification of Associated Interactions. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 1470-1480.	5.4	13
2	Pairing interactions between nucleobases and ligands in aptamer:ligand complexes of riboswitches: crystal structure analysis, classification, optimal structures, and accurate interaction energies. <i>Rna</i> , 2019, 25, 1274-1290.	3.5	8
3	Estimating Strengths of Individual Hydrogen Bonds in RNA Base Pairs: Toward a Consensus between Different Computational Approaches. <i>ACS Omega</i> , 2019, 4, 7354-7368.	3.5	21
4	Unfolding Transitions of Peripheral Subunit Binding Domains Show Cooperative Behavior. <i>Journal of Physical Chemistry B</i> , 2019, 123, 3441-3451.	2.6	0
5	Going beyond base-pairs: topology-based characterization of base-multiplets in RNA. <i>Rna</i> , 2019, 25, 573-589.	3.5	10
6	Evidence for Hidden Involvement of N3-Protonated Guanine in RNA Structure and Function. <i>ACS Omega</i> , 2019, 4, 699-709.	3.5	8
7	Consequences of Mg ²⁺ binding on the geometry and stability of RNA base pairs. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 21934-21948.	2.8	20
8	How Does Mg ²⁺ Modulate the RNA Folding Mechanism: A Case Study of the G:C W:W Trans Basepair. <i>Biophysical Journal</i> , 2017, 113, 277-289.	0.5	12
9	Structural landscape of base pairs containing post-transcriptional modifications in RNA. <i>Rna</i> , 2017, 23, 847-859.	3.5	29
10	Higher order structures involving post transcriptionally modified nucleobases in RNA. <i>RSC Advances</i> , 2017, 7, 35694-35703.	3.6	9
11	Effects of point mutations on the thermostability of <i>B. subtilis</i> lipase: investigating nonadditivity. <i>Journal of Computer-Aided Molecular Design</i> , 2016, 30, 899-916.	2.9	11
12	RNABP COGEST: a resource for investigating functional RNAs. <i>Database: the Journal of Biological Databases and Curation</i> , 2015, 2015, .	3.0	15
13	The role of N7 protonation of guanine in determining the structure, stability and function of RNA base pairs. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 26249-26263.	2.8	27
14	Understanding the Thermostability and Activity of <i>Bacillus subtilis</i> Lipase Mutants: Insights from Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2015, 119, 392-409.	2.6	58
15	Feasibility of occurrence of different types of protonated base pairs in RNA: a quantum chemical study. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 18383-18396.	2.8	36
16	Why Does Substitution of Thymine by 6-Ethynylpyridone Increase the Thermostability of DNA Double Helices?. <i>Journal of Physical Chemistry B</i> , 2014, 118, 6586-6596.	2.6	17
17	Protonation of Base Pairs in RNA: Context Analysis and Quantum Chemical Investigations of Their Geometries and Stabilities. <i>Journal of Physical Chemistry B</i> , 2011, 115, 1469-1484.	2.6	44
18	On the role of Hoogsteen:Hoogsteen interactions in RNA: Ab initio investigations of structures and energies. <i>Rna</i> , 2010, 16, 942-957.	3.5	39

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19	On the Role of the cis Hoogsteen/Sugar-Edge Family of Base Pairs in Platforms and Tripletsâ€”Quantum Chemical Insights into RNA Structural Biology. <i>Journal of Physical Chemistry B</i> , 2010, 114, 3307-3320.	2.6	33
20	MD simulations of ligand-bound and ligand-free aptamer: Molecular level insights into the binding and switching mechanism of the <i>add</i> A-riboswitch. <i>Rna</i> , 2009, 15, 1673-1692.	3.5	66
21	Modeling the noncovalent interactions at the metabolite binding site in purine riboswitches. <i>Journal of Molecular Modeling</i> , 2009, 15, 633-649.	1.8	27
22	Comparative modeling of thioredoxin glutathione reductase from <i>Schistosoma mansoni</i> : A multifunctional target for antischistosomal therapy. <i>Journal of Molecular Graphics and Modelling</i> , 2009, 27, 665-675.	2.4	24
23	Trans Hoogsteen/Sugar Edge Base Pairing in RNA. Structures, Energies, and Stabilities from Quantum Chemical Calculations. <i>Journal of Physical Chemistry B</i> , 2009, 113, 1743-1755.	2.6	55
24	A Theoretical Study on Interaction of Small Gold Clusters Au _n (n = 4, 6, 8) with xDNA Base Pairs. <i>Journal of Biomolecular Structure and Dynamics</i> , 2009, 27, 65-81.	3.5	20
25	Quantum Chemical Studies of Structures and Binding in Noncanonical RNA Base pairs: The Trans Watson-Crick/Watson-Crick Family. <i>Journal of Biomolecular Structure and Dynamics</i> , 2008, 25, 709-732.	3.5	50
26	Noncanonical Base Pairing in RNA: Topological and NBO Analysis of Hoogsteen Edge - Sugar Edge Interactions. <i>Lecture Notes in Computer Science</i> , 2008, , 379-386.	1.3	6
27	Theoretical analysis of noncanonical base pairing interactions in RNA molecules. <i>Journal of Biosciences</i> , 2007, 32, 809-825.	1.1	46
28	Base pairing in RNA structures: A computational analysis of structural aspects and interaction energies. <i>Journal of Chemical Sciences</i> , 2007, 119, 525-531.	1.5	24
29	Non-Canonical Base Pairs and Higher Order Structures in Nucleic Acids: Crystal Structure Database Analysis. <i>Journal of Biomolecular Structure and Dynamics</i> , 2006, 24, 149-161.	3.5	60
30	1,3-Dipolare Cycloadditionen, 92 Reaktionen aliphatischer Diazoverbindungen mit vierfach Acceptorsubstituierten Ethylenen. <i>Chemische Berichte</i> , 1987, 120, 153-158.	0.2	10
31	1,3-Dipolare Cycloadditionen, 93 Ãœberraschungen bei der Umsetzung des 2,3-DicyanfumarsÃ¤ure-dimethylesters mit Diazomethan. <i>Chemische Berichte</i> , 1987, 120, 159-169.	0.2	21
32	The Astounding Reaction of Diazomethane with Dimethyl 2,3-Dicyanofumarate. <i>Heterocycles</i> , 1986, 24, 2429.	0.7	7
33	Thermal and photochemical transformations of 1-(aryloxy)-n-arylidene-2-naphthylamines. <i>Journal of Organic Chemistry</i> , 1980, 45, 3182-3186.	3.2	6